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SELECTING PROCEDURES FOR OPTIMAL SUBSET OF REGRESSION VARIABLES*

by

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Recently, a number of methods have been developed for selecting the "best" or at least a "good" subset of variables in regression analysis. For various reasons, we may be interested in including only a subset, say, of size r < p, the number of independent variables. Various authors have considered this problem and a variety of techniques are presently being used to construct such subsets.

Arvesen and McCabe (1975) proposed a procedure for selecting a subset within a class of subsets with t (fixed) independent variables, taking into account the statistical variation of the residual mean squares. Huang and Panchapakesan (1982) proposed a selection procedure based on the expected residual sums of squares. Hsu and Huang (1982) studied a sequential selection procedure for good regression models.

In this paper, we are interested in deriving an optimal decision procedure based on residual mean squares to select a subset excluding all "inferior" independent variables. This kind of optimality criterion is related to the approach of Gupta and Huang (1977).

Let $\pi_0, \pi_1, \ldots, \pi_k$ denote k+1 normal populations with unknown variances $\sigma_0^2, \sigma_1^2, \ldots, \sigma_k^2$ respectively. Assume that σ_0^2 is known. A population (model) is said to be superior (or good) if $\sigma_1^2 < \Delta \sigma_0^2$, to be inferior (or bad) if

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 $\sigma_i^2 \ge \Delta \ \sigma_0^2$, where Δ is a specified constant greater than 1. Let Ω be the parameter space which is the collection of all possible parameters.

Let CD stand for a correct decision which is defined to be the selection of any subset which excludes all the inferior populations.

Assuming the following model

$$(1) Y = X\underline{\beta} + \epsilon$$

where $X = \begin{bmatrix} 1 \\ 1 \end{bmatrix}, X_1, \dots, X_{p-1} \end{bmatrix}$ is an nxp known matrix of rank $p \le n$, $\underline{\beta}' = (\beta_0, \beta_1, \dots, \beta_{p-1})$ is a 1xp parameter vector, and $\underline{\xi} \sim N(\underline{0}, \sigma_0^2 I_n)$, and $\underline{I}' = \begin{bmatrix} 1 \\ 1 \end{bmatrix}, \dots, 1 \end{bmatrix}_{1 \times N}$, I_n is an identity matrix with nxn.

In what follows, (1) which has p-1 independent variables, will be viewed as the true model. Without loss of generality we can assume that σ_0^2 = 1. Consider the models for any r, $2 \le r \le p-1$,

(2)
$$\underline{Y} = X_{ri} \underline{\beta}_{ri} + \underline{\epsilon}_{ri}$$

where X_{ri} is an nxr matrix of rank r with $X_{11}^t = [1, \dots, 1]_{1xn}$, β_{ri} is a rxl parameter vector, and $\xi_{ri} \sim N(\underline{0}, \sigma_{ri}^2|_{1})$, $i = 1, 2, \dots, k_r \sim \binom{p+1}{r-1}$. Let

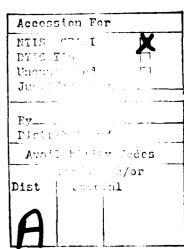
 $k = \sum_{r=2}^{p-1} k_r$. It should be noted that in stating the reduced model (2), our comparisons of models are made under the true model assumptions. The goal is to include all the designs X_{ri} (or sets of independent variables) associated with $\sigma_{[j]}^2$, $j = 1, \ldots, k-t$, where $\sigma_{[1]}^2 \leq \sigma_{[2]}^2 \leq \ldots \leq \sigma_{[k-t]}^2$ are ordered values from some of σ_{ri} 's, $i = 1, \ldots, k_r$, $r = 2, \ldots, p-1$.

Note that for any r, $2 \le r \le p-1$, if

$$SS_{ri} = Y' \{1 - X_{ri} (X_{ri} X_{ri})^{-1} X_{ri} \} Y = Y' Q_{i} Y,$$

then





$$ss_{ri} \sim \chi^2 \{v_r, (xg) \cdot Q_{ri}(xg)/2\}$$

(under the true model), where v_r = n-r, for $1 \le i \le k_r$. Note that the noncentrality parameter, in general, is not zero, and that

$$\sigma_{ri}^2 = 1 + (X\underline{\beta})'Q_{ri}(X\underline{\beta})/v_r.$$

Now we need some notation. Deleting a set of β_i 's without specifying which ones are deleted, we use ri to denote the special subset that is not deleted. For example, if p=3, r=2 then there are three subsets with size 2; namely, $\{\beta_1,\beta_2\}$, $\{\beta_1,\beta_3\}$ and $\{\beta_2,\beta_3\}$. Then rl denotes the set $\{\beta_1,\beta_2\}$, r2 denotes $\{\beta_2,\beta_3\}$ and r3 denotes $\{\beta_1,\beta_3\}$. Then, we use $\tilde{\beta}$ to denote the vector with the following subsets: $\{\beta_1,\beta_2\}$ with $(\beta_1,\beta_2,0)$, $\{\beta_1,\beta_3\}$ with $(\beta_1,0,\beta_3)$, and $\{\beta_2,\beta_3\}$ with $(0,\beta_2,\beta_3)$, where 0 is the parameter value which is omitted from the true model of the appropriate β_i 's. Thus, in the following, we will use $\Omega_{0,\mathbf{r}1}$ to denote those sets of $\tilde{\beta}$ as described above with the further condition that $\sigma_{\mathbf{r}i}^2=\sigma_0^2=1$. Similarly, $\Omega_{1,\mathbf{r}i}$ will be used to denote the sets of $\tilde{\beta}$ as described above with the further restriction that $\sigma_{\mathbf{r}i}^2\geq\Delta$. Formally, we write

$$\Omega_{0,ri} = \{\tilde{g} | \sigma_{ri}^2 = 1\},$$

and

$$\Omega_{1,ri} = \{\tilde{\beta} | \sigma_{ri}^2 \geq \Delta\},$$

where $i = 1, ..., k_r$; r = 2, ..., p-1, and let

$$\Omega_1 = \begin{array}{c} p-1 & k \\ \Gamma & \Gamma \\ \Gamma & \Gamma & \Gamma \\ \Gamma & \Gamma & \Gamma \end{array}$$

$$\Omega_0 = \bigcap_{r=2}^{p-1} \bigcap_{i=1}^{k} \Omega_{0,ri}.$$

Let $g_{\sigma_{ri}^2}(s_{ri})$ denote the probability density of s_{ri} depending on the

parameter σ_{ri}^2 , where $S_{ri} = \frac{SS_{ri}}{v_r}$, $i = 1,...,k_r$; r = 2,...,p-1.

Consider a family of hypotheses testing problems as follows:

(3)
$$H_{0,ri}: \tilde{\beta} \in \Omega_0 \text{ vs } K_{ri}: \tilde{\beta} \in \Omega_{1,ri};$$

i = 1,...,p-1, r = 2,...,p-1. A test of the hypotheses (3) will be defined to be a vector $(\phi_1(\underline{y}),\ldots,\phi_k(\underline{y}))$, where the elements of the vector are ordinary test functions; when \underline{y} is observed we reject H_0 , t with probability $\phi_t(\underline{y})$, $1 \le t \le k$. The power function of a test (ϕ_1,\ldots,ϕ_k) is defined to be the vector $(p_1(\underline{\tilde{\beta}}),\ldots,p_k(\underline{\tilde{\beta}}))$ where

$$p_{t}(\tilde{\beta}) = E_{\tilde{\beta}} \varphi_{t}(\tilde{Y}),$$

 $1 \leq t \leq k.$ Let $S(\gamma)$ be the set of all tests (ϕ_1, \ldots, ϕ_k) such that

(4)
$$E_{\tilde{\beta}} \varphi_{t}(\underline{Y}) \leq \gamma, \quad \tilde{\underline{\beta}} \in \Omega_{0}.$$

We define $\varphi^0 = (\varphi_1^0, \dots, \varphi_k^0)$ as

$$\varphi_{ri}^{0}(\underline{y}) = \begin{cases} 1, & \text{if } g_{\Delta}(s_{ri}) \geq c \ g_{1}(s_{ri}), \\ \\ 0, & \text{if } g_{\Delta}(s_{ri}) < c \ g_{1}(s_{ri}), \end{cases}$$

such that $E_{\tilde{\beta}}^{0} \varphi_{ri}^{0}(\underline{Y}) = \gamma$, $\tilde{\beta} \in \Omega_{0}$, where s_{ri} is the observed value of S_{ri} . It can be shown that φ^{0} maximizes

$$\min_{\substack{1 \le t \le k \\ \tilde{\beta} \in \Omega_{1,t}}} \inf_{\tilde{\beta} \in \Omega_{1,t}} E_{\tilde{\beta}} \varphi_{t}(\underline{Y})$$

among all tests ϕ = (ϕ_1,\ldots,ϕ_k) \in $S(\gamma)$ (cf. Gupta and Huang (1977)).

To determine the constant c, we proceed follows: for a given $\eta > 0$, there exists a smallest positive integer k_0 such that

$$\frac{a_{k_0}}{n} < 1$$
 and $\frac{a_{k_0+1}}{a_{k_0}} + \frac{a_{k_0}}{n} \le 1$,

where

$$a_{\ell}(s_{ri}) = \frac{e^{-\lambda}r_{\lambda_r^{\ell}}}{\ell!} \left[\frac{v_r s_{ri}}{2}\right]^{\ell} \frac{\Gamma(\frac{1}{2}v_r)}{\Gamma(\frac{1}{2}v_r + \ell)},$$

 $\ell = 0, 1, 2, ...; \lambda_r = \frac{(\Delta - 1)v_r}{2}$. For this k_0 , it can be shown that

$$0 < \frac{g_{\Delta}(s_{ri})}{g_{1}(s_{ri})} - \sum_{k=0}^{k_{0}-1} a_{k}(s_{ri}) = \sum_{k=0}^{\infty} a_{k_{0}+k} \leq \eta,$$

where

$$\frac{g_{\Delta}(s_{ri})}{g_{1}(s_{ri})} = \sum_{\ell=0}^{\infty} a_{\ell}.$$

Thus, approximately,

$$\frac{g_{\Delta}(s_{ri})}{g_{1}(s_{ri})} = \sum_{\ell=0}^{k-1} a_{\ell}(s_{ri})$$

with error less than n. For $\tilde{\beta} \in \Omega_0$,

$$E_{\tilde{\beta}} \varphi^{0}(Y) = P_{\tilde{\beta}} \{g_{\Delta}(S_{ri}) \ge c \ g_{1}(S_{ri})\}$$

$$= P_{\tilde{\beta}} \{\sum_{\ell=0}^{k_{0}-1} a_{\ell}(S_{ri}) \ge c\}$$

$$= \int_{0}^{\infty} I_{k_{0}-1} (s_{ri})g_{1}(s_{ri})ds_{ri} = \gamma,$$

$$\{\sum_{\ell=0}^{k_{0}-1} a_{\ell}(s_{ri}) \ge c\}$$

where $g_1(s_{ri})$ is the central χ^2 with v_r degrees of freedom and $l_{\Lambda}(x) = 0$ for $x \notin A$, $l_{\Lambda}(x) = 1$ for $x \in A$. The constant c can be determined.

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Recently, a number of methods have been developed for selecting the "best" or at least a "good" subset of variables in regression analysis. For various reasons, we may be interested in including only a subset, say, of size r < p, the number of independent variables. Various authors have considered this problem and a variety of techniques are presently being used to construct such subsets. In this paper, we are interested in deriving an optimal decision procedure based on residual mean squares to select a subset excluding all "inferior" independent variables. This kind of optimality criterion is related to the approach of Gupta and Huang (1977).

